

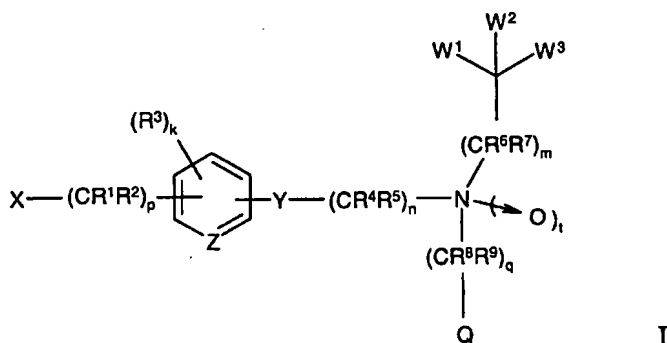
Amendment to the Claims:

This listing of claims will replace all prior versions and listing of claims in the application:

Listing of Claims:

1-48 (Cancelled).

49. (Currently amended): A compound of Formula I:



wherein:

X is selected from C₁-C₈ alkyl, halo, -OR¹⁰, -NR¹⁴R¹⁵, nitro, cyano, -COOR¹⁰, -COR¹³, -OCOR¹³, -N(R¹⁷)COR¹³, -N(R¹⁷)CONR¹⁴R¹⁵, -N(R¹⁷)COOR¹³, -SO₃H, -SO₂NR¹⁴R¹⁵, -C(=NR¹⁷)NR¹⁴R¹⁵, -N(R¹⁷)SO₂R¹⁶, and a 5 or 6-membered heterocyclic group;

or X and an adjacent R³, taken together with the atoms to which they are bonded, form an alkylenedioxy moiety;

Z is CH, CR³ or N, wherein when Z is CH or CR³, k is 0-4 and t is 0 or 1, and when Z is N, k is 0-3 and t is 0;

Y is selected from -O-, -S-, -N(R¹⁰)-, and -C(R⁴)(R⁵)-;

W¹ is selected from [[C₄-C₆ alkyl],] C₃-C₈ cycloalkyl, aryl and Het, wherein said [[C₄-C₆ alkyl],] C₃-C₈ cycloalkyl, aryl and Het are optionally unsubstituted or substituted with one or more groups independently selected from halo, cyano, nitro, C₁-C₆ alkyl, C₃-C₆ alkenyl, C₃-C₆ alkynyl, -C₀-C₆ alkyl-CO₂R¹⁰, -C₀-C₆ alkyl-C(O)SR¹⁰, -C₀-C₆ alkyl-CONR¹¹R¹², -C₀-C₆ alkyl-COR¹³, -C₀-C₆ alkyl-NR¹¹R¹², -C₀-C₆ alkyl-SR¹⁰, -C₀-C₆ alkyl-OR¹⁰, -C₀-C₆ alkyl-SO₃H, -C₀-C₆ alkyl-SO₂NR¹¹R¹², -C₀-C₆ alkyl-SO₂R¹⁰, -C₀-C₆ alkyl-SOR¹³, -C₀-C₆ alkyl-OCOR¹³, -C₀-C₆ alkyl-OC(O)NR¹¹R¹², -C₀-C₆ alkyl-OC(O)OR¹³, -C₀-C₆ alkyl-NR¹¹C(O)OR¹³, -C₀-C₆ alkyl-NR¹¹C(O)NR¹¹R¹², and

$-C_0-C_6$ alkyl- $NR^{11}COR^{13}$, where said C_1-C_6 alkyl, is optionally unsubstituted or substituted by one or more halo substituents;

W^2 is selected from H, halo, C_1-C_6 alkyl, C_2-C_6 alkenyl, C_2-C_6 alkynyl, $-C_0-C_6$ alkyl- $NR^{11}R^{12}$, $-C_0-C_6$ alkyl- SR^{10} , $-C_0-C_6$ alkyl- OR^{10} , $-C_0-C_6$ alkyl- CO_2R^{10} , $-C_0-C_6$ alkyl- $C(O)SR^{10}$, $-C_0-C_6$ alkyl- $CONR^{11}R^{12}$, $-C_0-C_6$ alkyl- COR^{13} , $-C_0-C_6$ alkyl- $OCOR^{13}$, $-C_0-C_6$ alkyl- $OCONR^{11}R^{12}$, $-C_0-C_6$ alkyl- $NR^{11}CONR^{11}R^{12}$, $-C_0-C_6$ alkyl- $NR^{11}COR^{13}$, $-C_0-C_6$ alkyl-Het, $-C_0-C_6$ alkyl-aryl and $-C_0-C_6$ alkyl- C_3-C_7 cycloalkyl, wherein said C_1-C_6 alkyl is optionally unsubstituted or substituted by one or more halo substituents, and wherein the C_3-C_7 cycloalkyl, aryl and Het moieties of said $-C_0-C_6$ alkyl-Het, $-C_0-C_6$ alkyl-aryl and $-C_0-C_6$ alkyl- C_3-C_7 cycloalkyl are optionally unsubstituted or substituted with one or more groups independently selected from halo, cyano, nitro, C_1-C_6 alkyl, C_3-C_6 alkenyl, C_3-C_6 alkynyl, $-C_0-C_6$ alkyl- CO_2R^{10} , $-C_0-C_6$ alkyl- $C(O)SR^{10}$, $-C_0-C_6$ alkyl- $CONR^{11}R^{12}$, $-C_0-C_6$ alkyl- COR^{13} , $-C_0-C_6$ alkyl- $NR^{11}R^{12}$, $-C_0-C_6$ alkyl- SR^{10} , $-C_0-C_6$ alkyl- OR^{10} , $-C_0-C_6$ alkyl- SO_3H , $-C_0-C_6$ alkyl- $SO_2NR^{11}R^{12}$, $-C_0-C_6$ alkyl- SO_2R^{10} , $-C_0-C_6$ alkyl- SOR^{13} , $-C_0-C_6$ alkyl- $OCOR^{13}$, $-C_0-C_6$ alkyl- $OC(O)NR^{11}R^{12}$, $-C_0-C_6$ alkyl- $OC(O)OR^{13}$, $-C_0-C_6$ alkyl- $NR^{11}C(O)OR^{13}$, $-C_0-C_6$ alkyl- $NR^{11}C(O)NR^{11}R^{12}$, and $-C_0-C_6$ alkyl- $NR^{11}COR^{13}$, where said C_1-C_6 alkyl, is optionally unsubstituted or substituted by one or more halo substituents;

W^3 is selected from the group consisting of: H, halo, C_1-C_6 alkyl, $-C_0-C_6$ alkyl- $NR^{11}R^{12}$, $-C_0-C_6$ alkyl- SR^{10} , $-C_0-C_6$ alkyl- OR^{10} , $-C_0-C_6$ alkyl- CO_2R^{10} , $-C_0-C_6$ alkyl- $C(O)SR^{10}$, $-C_0-C_6$ alkyl- $CONR^{11}R^{12}$, $-C_0-C_6$ alkyl- COR^{13} , $-C_0-C_6$ alkyl- $OCOR^{13}$, $-C_0-C_6$ alkyl- $OCONR^{11}R^{12}$, $-C_0-C_6$ alkyl- $NR^{11}CONR^{11}R^{12}$, $-C_0-C_6$ alkyl- $NR^{11}COR^{13}$, $-C_0-C_6$ alkyl-Het, $-C_1-C_6$ alkyl-aryl and $-C_1-C_6$ alkyl- C_3-C_7 cycloalkyl, wherein said C_1-C_6 alkyl is optionally unsubstituted or substituted by one or more halo substituents;

Q is selected from C_3-C_8 cycloalkyl, Ar and Het; wherein said C_3-C_8 cycloalkyl, Ar and Het are optionally unsubstituted or substituted with one or more groups independently selected from halo, cyano, nitro, C_1-C_6 alkyl, C_3-C_6 alkenyl, C_3-C_6 alkynyl, $-C_0-C_6$ alkyl- CO_2R^{10} , $-C_0-C_6$ alkyl- $C(O)SR^{10}$, $-C_0-C_6$ alkyl- $CONR^{11}R^{12}$, $-C_0-C_6$ alkyl- COR^{13} , $-C_0-C_6$ alkyl- $NR^{11}R^{12}$, $-C_0-C_6$ alkyl- SR^{10} , $-C_0-C_6$ alkyl- OR^{10} , $-C_0-C_6$ alkyl- SO_3H , $-C_0-C_6$ alkyl- $SO_2NR^{11}R^{12}$, $-C_0-C_6$ alkyl- SO_2R^{10} , $-C_0-C_6$ alkyl- SOR^{13} , $-C_0-C_6$ alkyl- $OCOR^{13}$, $-C_0-C_6$ alkyl- $OC(O)NR^{11}R^{12}$, $-C_0-C_6$ alkyl- $OC(O)OR^{13}$, $-C_0-C_6$ alkyl- $NR^{11}C(O)OR^{13}$,

$-C_0-C_6$ alkyl- $NR^{11}C(O)NR^{11}R^{12}$, and $-C_0-C_6$ alkyl- $NR^{11}COR^{13}$, where said C_1-C_6 alkyl is optionally unsubstituted or substituted by one or more halo substituents;

p is 0-8;

n is 3;

m is 0 or 1;

q is 0 or 1;

t is 0 or 1;

each R^1 and R^2 are independently selected from H, halo, C_1-C_6 alkyl, C_3-C_6 alkenyl, C_3-C_6 alkynyl, $-C_0-C_6$ alkyl- $NR^{11}R^{12}$, $-C_0-C_6$ alkyl- OR^{10} , $-C_0-C_6$ alkyl- SR^{10} , $-C_1-C_6$ alkyl-Het, $-C_1-C_6$ alkyl-Ar and $-C_1-C_6$ alkyl- C_3-C_7 cycloalkyl, or R^1 and R^2 together with the carbon to which they are attached form a 3-5 membered carbocyclic or heterocyclic ring, wherein said heterocyclic ring contains one, or more heteroatoms selected from N, O, and S, where any of said C_1-C_6 alkyl is optionally unsubstituted or substituted by one or more halo substituents;

each R^3 is the same or different and is independently selected from halo, cyano, nitro, C_1-C_6 alkyl, C_3-C_6 alkenyl, C_3-C_6 alkynyl, $-C_0-C_6$ alkyl-Ar, $-C_0-C_6$ alkyl-Het, $-C_0-C_6$ alkyl- C_3-C_7 cycloalkyl, $-C_0-C_6$ alkyl- CO_2R^{10} , $-C_0-C_6$ alkyl- $C(O)SR^{10}$, $-C_0-C_6$ alkyl- $CONR^{11}R^{12}$, $-C_0-C_6$ alkyl- COR^{13} , $-C_0-C_6$ alkyl- $NR^{11}R^{12}$, $-C_0-C_6$ alkyl- SR^{10} , $-C_0-C_6$ alkyl- OR^{10} , $-C_0-C_6$ alkyl- SO_3H , $-C_0-C_6$ alkyl- $SO_2NR^{11}R^{12}$, $-C_0-C_6$ alkyl- SO_2R^{10} , $-C_0-C_6$ alkyl- SOR^{13} , $-C_0-C_6$ alkyl- $OCOR^{13}$, $-C_0-C_6$ alkyl- $OC(O)NR^{11}R^{12}$, $-C_0-C_6$ alkyl- $OC(O)OR^{13}$, $-C_0-C_6$ alkyl- $NR^{11}C(O)OR^{13}$, $-C_0-C_6$ alkyl- $NR^{11}C(O)NR^{11}R^{12}$, and $-C_0-C_6$ alkyl- $NR^{11}COR^{13}$, wherein said C_1-C_6 alkyl is optionally unsubstituted or substituted by one or more halo substituents;

each R^4 and R^5 is independently selected from H, halo, C_1-C_6 alkyl, $-C_0-C_6$ alkyl-Het, $-C_0-C_6$ alkyl-Ar and $-C_0-C_6$ alkyl- C_3-C_7 cycloalkyl;

R^6 and R^7 are each independently selected from H, halo, C_1-C_6 alkyl, $-C_0-C_6$ alkyl-Het, $-C_0-C_6$ alkyl-Ar and $-C_0-C_6$ alkyl- C_3-C_7 cycloalkyl;

R^8 and R^9 are each independently selected from H, halo, C_1-C_6 alkyl, $-C_0-C_6$ alkyl-Het, $-C_0-C_6$ alkyl-Ar and $-C_0-C_6$ alkyl- C_3-C_7 cycloalkyl;

R^{10} is selected from H, C_1-C_6 alkyl, C_3-C_6 alkenyl, C_3-C_6 alkynyl, $-C_0-C_6$ alkyl-Ar, $-C_0-C_6$ alkyl-Het and $-C_0-C_6$ alkyl- C_3-C_7 cycloalkyl;

each R^{11} and each R^{12} are independently selected from H, C_1-C_6 alkyl, C_3-C_6 alkenyl, C_3-C_6 alkynyl, $-C_0-C_6$ alkyl-Ar, $-C_0-C_6$ alkyl-Het and $-C_0-C_6$ alkyl- C_3-C_7 cycloalkyl, or R^{11}

and R¹² together with the nitrogen to which they are attached form a 4-7 membered heterocyclic ring which optionally contains one or more additional heteroatoms selected from N, O, and S;

R¹³ is selected from C₁-C₆ alkyl, C₃-C₆ alkenyl, C₃-C₆ alkynyl, -C₀-C₆ alkyl-Ar, -C₀-C₆ alkyl-Het and -C₀-C₆ alkyl-C₃-C₇ cycloalkyl;

R¹⁴ and R¹⁵ are each independently selected from H, C₁-C₆ alkyl, C₃-C₆ alkenyl, C₃-C₆ alkynyl, -C₀-C₆ alkyl-Ar, -C₀-C₆ alkyl-Het, -C₀-C₆ alkyl-C₃-C₇ cycloalkyl, -C₀-C₆ alkyl-O-Ar, -C₀-C₆ alkyl-O-Het, -C₀-C₆ alkyl-O-C₃-C₇ cycloalkyl, -C₀-C₆ alkyl-S(O)_x-C₁-C₆ alkyl, -C₀-C₆ alkyl-S(O)_x-Ar, -C₀-C₆ alkyl-S(O)_x-Het, -C₀-C₆ alkyl-S(O)_x-C₃-C₇ cycloalkyl, -C₀-C₆ alkyl-NH-Het, -C₀-C₆ alkyl-NH-C₃-C₇ cycloalkyl, -C₀-C₆ alkyl-N(C₁-C₄ alkyl)-Ar, -C₀-C₆ alkyl-N(C₁-C₄ alkyl)-Het, -C₀-C₆ alkyl-N(C₁-C₄ alkyl)-C₃-C₇ cycloalkyl, -C₀-C₆ alkyl-Ar, -C₀-C₆ alkyl-Het and -C₀-C₆ alkyl-C₃-C₇ cycloalkyl, where x is 0, 1 or 2, or R¹⁴ and R¹⁵, together with the nitrogen to which they are attached, form a 4-7 membered heterocyclic ring which optionally contains one or more additional heteroatoms selected from N, O, and S, wherein said C₁-C₆ alkyl is optionally substituted by one or more of the substituents independently selected from the group halo, -OH, -SH, -NH₂, -NH(unsubstituted C₁-C₆ alkyl), -N(unsubstituted C₁-C₆ alkyl)(unsubstituted C₁-C₆ alkyl), unsubstituted -OC₁-C₆ alkyl, -CO₂H, -CO₂(unsubstituted C₁-C₆ alkyl), -CONH₂, -CONH(unsubstituted C₁-C₆ alkyl), -CON(unsubstituted C₁-C₆ alkyl)(unsubstituted C₁-C₆ alkyl), -SO₃H, -SO₂NH₂, -SO₂NH(unsubstituted C₁-C₆ alkyl) and -SO₂N(unsubstituted C₁-C₆ alkyl)(unsubstituted C₁-C₆ alkyl);

R¹⁶ is C₁-C₆ alkyl, -C₀-C₆ alkyl-Ar or -C₀-C₆ alkyl-Het; and

R¹⁷ is H, C₁-C₆ alkyl, -C₀-C₆ alkyl-Ar or -C₀-C₆ alkyl-Het;

wherein each Ar or aryl independently represent a substituted or unsubstituted carbocyclic aromatic group, which may be optionally fused to another carbocyclic aromatic group moiety or to a cycloalkyl group moiety, wherein said Ar or aryl is optionally substituted by one or more of the substituents independently selected from the group halo, cyano, C₁-C₆ alkyl, C₁-C₆ haloalkyl, -C₀-C₆ alkyl-OH, -C₀-C₆ alkyl-SH, -C₀-C₆ alkyl-NR'R", C₃-C₆ alkenyl, -OC₁-C₆alkyl, -OC₁-C₆ alkenyl, -C₀-C₆ alkyl-COR', -C₀-C₆ alkyl-CO₂R', -C₀-C₆ alkyl-CONR'R", -OC₀-C₆ alkyl-CO₂H, -OC₂-C₆ alkyl-NR'R",

-C₀-C₆ alkyl-C(=NR')NR'R", and -C₀-C₆ alkyl-SO₂NR'R", wherein each R' and R" are independently selected from H and unsubstituted C₁-C₆ alkyl,

each Het independently represents a monocyclic 5- to 7-membered, a bicyclic 7- to 10-membered or ~~an 11- to 18-membered trieyclie~~ a tricyclic 11- to 18-membered heterocyclic ring group which is saturated, unsaturated or aromatic, and consists of carbon atoms and from one to three heteroatoms selected from N, O and S, wherein the N or S heteroatoms of said Het are optionally oxidized or the N heteroatom is optionally quaternized, wherein said Het is optionally unsubstituted or substituted by one or more of the substituents independently selected from the group halo, cyano, C₁-C₆ alkyl (which specifically includes C₁-C₆ haloalkyl, -C₀-C₆ alkyl-OH, -C₀-C₆ alkyl-SH and -C₀-C₆ alkyl-NR'R"), C₃-C₆ alkenyl, oxo, -OC₁-C₆alkyl, -OC₁-C₆ alkenyl, -C₀-C₆ alkyl-COR', -C₀-C₆ alkyl-CO₂R', -C₀-C₆ alkyl-CONR'R", -OC₀-C₆ alkyl-CO₂H, -OC₂-C₆ alkyl-NR'R", -C₀-C₆ alkyl-C(=NR')NR'R" and -C₀-C₆ alkyl-SO₂NR'R", wherein each R' and R" are independently selected from H and unsubstituted C₁-C₆ alkyl;

provided that X is not COOR¹⁰ when Y is -O-, p is 0-8, n is 3, m is 1, q is 0 or 1, t is 0, each R¹ and R² is independently selected from H, C₁-C₆ alkyl, -OH, -O-C₁-C₆ alkyl, -SH, and -S-C₁-C₆ alkyl, each R⁴, R⁵, R⁶, R⁷, R⁸ and R⁹ are independently H or C₁-C₄ alkyl, k is 0 or 1, W³ is H, W¹ and W² are each independently selected from C₃-C₈ cycloalkyl and aryl and R³ and Q are as defined above; or

provided that the compound is not

5-[3-[(3,4-dichlorophenyl)methyl][2-(2-naphthalenyl)ethyl]amino]propoxy]-3-methoxy-1,2-benzenedicarboxylic acid, or

5-[3-[(3,4-dichlorophenyl)methyl][2-(2-naphthalenyl)ethyl]amino]propoxy]-3-methoxy-1,2-benzenedicarboxylic acid, dimethyl ester [[.]];

4-[[[2-(4-carboxyphenoxy)ethyl][2-[2-[(5-phenylpentyl)oxy]phenyl]ethyl]amino]methyl] benzoic acid;

4-[[[2-[4-(ethoxycarbonyl)phenoxy]ethyl][2-[2-(octyloxy)phenyl]ethyl]amino]methyl] benzoic acid methyl ester;

4-[[[2-(4-carboxyphenoxy)ethyl][2-[2-(octyloxy)phenyl]ethyl]amino]methyl] benzoic acid;

α -[[[3-(4-fluorophenyl)-1,1-dimethylpropyl](phenylmethyl)amino]methyl]-3-(phenylmethoxy) benzenemethanol hydrochloride;

~~N-[2-(4-amino-3,5-dichlorophenyl)ethyl]-4-fluoro-N-(phenylmethyl)-benzenepropanamine monohydrochloride;~~

~~N-[2-(4-amino-3,5-dichlorophenyl)ethyl]-4-chloro-N-(phenylmethyl)-benzenepropanamine monohydrochloride;~~

~~4-amino-3,5-dichloro- α -[[[3-(4-fluorophenyl)propyl](phenylmethyl)amino]methyl]-benzenemethanol monohydrochloride;~~

~~4-amino-3,5-dichloro- α -[[[3-(4-chlorophenyl)propyl](phenylmethyl)amino]methyl]-benzenemethanol monohydrochloride;~~

~~2-chloro-5-[2-[[3-(4-fluorophenyl)-1-methylpropyl](phenylmethyl)amino]-1-hydroxyethyl]-benzamide monohydrochloride;~~

~~4-[2-[[2-hydroxy-2-[4-(phenylmethoxy)phenyl]ethyl](phenylmethyl)amino]ethoxy]-benzeneacetamide;~~

~~4-[2-[[2-[3,4-bis(phenylmethoxy)phenyl]ethyl](phenylmethyl)amino]ethoxy]-benzenesulfonamide monohydrochloride;~~

~~(R)-3-(phenylmethoxy)- α -[[[3-(3-(phenylmethoxy)phenyl)propyl](phenylmethyl)amino]methyl]-benzenemethanol;~~

~~2,2-dichloro-acetic acid (R)-[benzyl [3-(3-benzyloxy-phenyl)-propyl]-amino]-(3-benzyloxy-phenyl)-ethyl ester;~~

~~3-amino- α -[[[3-(3,4-dimethoxyphenyl)-1-methylpropyl](phenylmethyl)amino]methyl]-4-(phenylmethoxy)-benzenemethanol;~~

~~α -[[[3-(3,4-dimethoxyphenyl)-1-methylpropyl](phenylmethyl)amino]methyl]-3-nitro-4-(phenylmethoxy)-benzenemethanol;~~

~~α -[[[3-(3,4-dimethoxyphenyl)-1-methylpropyl](phenylmethyl)amino]methyl]-3-nitro-5-(phenylmethoxy)-benzenemethanol;~~

~~3-amino- α -[[[3-(3,4-dimethoxyphenyl)-1-methylpropyl](phenylmethyl)amino]methyl]-5-(phenylmethoxy)-benzenemethanol; or~~

~~4-[2-[[2-(4-fluorophenoxy)ethyl](phenylmethyl)amino]ethyl]-1-piperazineacetic acid ethyl ester;~~

or a pharmaceutically acceptable salt or hydrate thereof.

50. (Previously presented): The compound according to claim 49, wherein p is 0, 1 or 2.

51. (Previously presented): The compound according to claim 49, wherein t is 0.

52. (Previously presented): The compound according to claim 49, wherein R¹ and R² are independently H or C₁-C₄ alkyl or R¹ and R² together with the carbon to which they are attached form a 3-5 membered carbocyclic ring.

53. (Previously presented): The compound according to claim 49, wherein k is 0 or 1.

54. (Previously presented): The compound according to claim 49, wherein R³ is selected from halo, C₁-C₄ alkyl and C₁-C₄ alkoxy.

55. (Previously presented): The compound according to claim 49, wherein X is selected from C₁-C₆ alkyl, halo, -OR¹⁰, -NR¹⁴R¹⁵, cyano, -COR¹³, -COOR¹⁰, -OCOR¹³, -N(R¹⁷)CONR¹⁴R¹⁵, -N(R¹⁷)COR¹³, -SO₂NR¹⁴R¹⁵, -N(R¹⁷)SO₂R¹⁶, and a 5 or 6-membered heterocyclic group or X and an adjacent R³, taken together with the atoms to which they are bonded, form an alkylenedioxy moiety.

56. (Previously presented): The compound according to claim 55, wherein R¹⁰ is H, C₁-C₄ alkyl or phenyl; R¹³ is H, C₁-C₄ alkyl, -C₀-C₄ alkyl-C₃-C₇ cycloalkyl, or -C₀-C₄ alkyl-phenyl; R¹⁴ and R¹⁵ are each independently selected from H, C₁-C₆ alkyl, -C₀-C₄ alkyl-Ar, -C₀-C₄ alkyl-Het, -C₀-C₄ alkyl-C₃-C₇ cycloalkyl, -C₀-C₄ alkyl-O-Ar, -C₀-C₄ alkyl-O-Het, -C₀-C₄ alkyl-O-C₃-C₇ cycloalkyl, -C₀-C₄ alkyl-S(O)₂-C₁-C₄ alkyl, -C₀-C₄ alkyl-S(O)₂-Ar, -C₀-C₄ alkyl-S(O)₂-Het, -C₀-C₄ alkyl-S(O)₂-C₃-C₇ cycloalkyl, -C₀-C₄ alkyl-NH-Ar, -C₀-C₄ alkyl-NH-Het, -C₀-C₄ alkyl-NH-C₃-C₇ cycloalkyl, -C₀-C₄ alkyl-N(C₁-C₄ alkyl)-Ar, -C₀-C₄ alkyl-N(C₁-C₄ alkyl)-Het, -C₀-C₄ alkyl-N(C₁-C₄ alkyl)-C₃-C₇ cycloalkyl, -C₀-C₄ alkyl-Ar, -C₀-C₄ alkyl-Het and -C₀-C₄ alkyl-C₃-C₇ cycloalkyl, or R¹⁴ and R¹⁵, together with the nitrogen to which they are attached, form a 4-7 membered heterocyclic ring which optionally contains one or more additional heteroatoms selected from N, O, and S, wherein said C₁-C₆ alkyl is optionally substituted by one or more of the substituents independently selected from the group halo,

-OH, -SH, -NH₂, -NH(unsubstituted C₁-C₄ alkyl), -N(unsubstituted C₁-C₄ alkyl)(unsubstituted C₁-C₄ alkyl), unsubstituted -OC₁-C₄ alkyl, -CO₂H, -CO₂(unsubstituted C₁-C₄ alkyl), -CONH₂, -CONH(unsubstituted C₁-C₄ alkyl), -CON(unsubstituted C₁-C₄ alkyl)(unsubstituted C₁-C₄ alkyl), -SO₃H, -SO₂NH₂, -SO₂NH(unsubstituted C₁-C₄ alkyl) and -SO₂N(unsubstituted C₁-C₄ alkyl)(unsubstituted C₁-C₄ alkyl); R¹⁶ is C₁-C₄ alkyl or phenyl; and R¹⁷ is H or C₁-C₄ alkyl.

57. (Previously presented): The compound according to claim 49 wherein each R⁴ and R⁵ are independently selected from H and C₁-C₃ alkyl.

Claim 58 (Cancelled).

59. (Previously presented): The compound according to claim 49, wherein R⁸ and R⁹ are each H.

60. (Previously presented): The compound according to claim 49, wherein Q is a substituted or unsubstituted phenyl or furanyl group or a benzo[1,3]dioxyl or benzo[1,4]dioxyl group containing one, two or three substituents selected from halo, C₁-C₄ alkyl; C₁-C₄ alkylthio; or -NR^{Q1}R^{Q2}, where R^{Q1} and R^{Q2} taken together with the nitrogen to which they are attached form a 4-7 membered heterocyclic ring, which may optionally contain one or more additional heteroatoms selected from N, O and S.

61. (Previously presented): The compound according to claim 60, wherein said substituents are selected from fluoro, chloro, trifluoromethyl, tert-butyl, isopropyl, methylthio and piperidin-1-yl.

62. (Previously presented): The compound according to claim 49, wherein m is 0 or m is 1 and R⁶ and R⁷ are each H.

63. (Previously presented): The compound according to claim 49, wherein W¹ is phenyl, naphthyl, thienyl, pyridyl, furanyl, pyrrolyl, cyclohexyl, cyclopentyl, morpholinyl, or pyrrolidinyl, where each phenyl, naphthyl, thienyl, pyridyl, furanyl, pyrrolyl, cyclohexyl,

cyclopentyl, morpholinyl, or pyrrolidinyl may be optionally substituted from 1 to 3 times with one or more of the substituents independently selected from C₁-C₄ alkyl, -OH, halo, -O-C₁-C₄ alkyl, and -C₁-C₄ haloalkyl.

64. (Previously presented): The compound according to claim 49, wherein W² is C₁-C₄ alkyl, C₂-C₄ alkynyl, C₃-C₆ cycloalkyl, aryl, Het hydroxy, aryloxy-, C₁-C₄ alkoxy-, -OCOC₁-C₄ alkyl, -OCOaryl, or -NR^{W1}R^{W2}, where R^{W1} and R^{W2} are independently H or C₁-C₄ alkyl or taken together with the nitrogen to which they are attached form a 4-7 membered heterocyclic ring, which may optionally contain one or more additional heteroatoms selected from N, O and S.

65. (Withdrawn): The compound according to claim 49, wherein W³ is H or C₁-C₄ alkyl.

66. (Currently amended): The compound according to claim 49, wherein X is selected from C₁-C₆ alkyl, halo, -OR¹⁰, -NR¹⁴R¹⁵, cyano, -COR¹³, -COOR¹⁰, -OCOR¹³, -N(R¹⁷)CONR¹⁴R¹⁵, -N(R¹⁷)COR¹³, -SO₂NR¹⁴R¹⁵, -N(R¹⁷)SO₂R¹⁶, and a 5 or 6-membered heterocyclic group or X and an adjacent R³, taken together with the atoms to which they are bonded, form an alkylenedioxy moiety, where R¹⁰ is H, C₁-C₄ alkyl or phenyl, R¹³ is H, C₁-C₄ alkyl, -C₀-C₄ alkyl-C₃-C₇ cycloalkyl, or -C₀-C₄ alkyl-phenyl, R¹⁴ and R¹⁵ are each independently selected from H, C₁-C₆ alkyl, -C₀-C₄ alkyl-Ar, -C₀-C₄ alkyl-Het, -C₀-C₄ alkyl-C₃-C₇ cycloalkyl, -C₀-C₄ alkyl-O-Ar, -C₀-C₄ alkyl-O-Het, -C₀-C₄ alkyl-O-C₃-C₇ cycloalkyl, -C₀-C₄ alkyl-S(O)₂-C₁-C₄ alkyl, -C₀-C₄ alkyl-S(O)₂-Ar, -C₀-C₄ alkyl-S(O)₂-Het, -C₀-C₄ alkyl-S(O)₂-C₃-C₇ cycloalkyl, -C₀-C₄ alkyl-NH-Ar, -C₀-C₄ alkyl-NH-Het, -C₀-C₄ alkyl-NH-C₃-C₇ cycloalkyl, -C₀-C₄ alkyl-N(C₁-C₄ alkyl)-Ar, -C₀-C₄ alkyl-N(C₁-C₄ alkyl)-Het, -C₀-C₄ alkyl-N(C₁-C₄ alkyl)-C₃-C₇ cycloalkyl, -C₀-C₄ alkyl-Ar, -C₀-C₄ alkyl-Het and -C₀-C₄ alkyl-C₃-C₇ cycloalkyl, or R¹⁴ and R¹⁵, together with the nitrogen to which they are attached, form a 4-7 membered heterocyclic ring which optionally contains one or more additional heteroatoms selected from N, O, and S, wherein said C₁-C₆ alkyl is optionally substituted by one or more of the substituents independently selected from the group halo, -OH, -SH, -NH₂, -NH(unsubstituted C₁-C₄ alkyl), -N(unsubstituted C₁-C₄ alkyl)(unsubstituted C₁-C₄ alkyl), unsubstituted -OC₁-C₄ alkyl,

-CO₂H, -CO₂(unsubstituted C₁-C₄ alkyl), -CONH₂, -CONH(unsubstituted C₁-C₄ alkyl),
-CON(unsubstituted C₁-C₄ alkyl)(unsubstituted C₁-C₄ alkyl), -SO₃H, -SO₂NH₂,
-SO₂NH(unsubstituted C₁-C₄ alkyl) and -SO₂N(unsubstituted C₁-C₄ alkyl)(unsubstituted
C₁-C₄ alkyl), R¹⁶ is C₁-C₄ alkyl or phenyl, and R¹⁷ is H or C₁-C₄ alkyl; p is 0, 1 or 2; R¹ and
R² are independently H or C₁-C₄ alkyl or R¹ and R² together with the carbon to which they are
attached form a 3-5 membered carbocyclic ring; k is 0 or k is 1 and R³ is halo, C₁-C₄ alkyl or
C₁-C₄ alkoxy; n is 3 and each R⁴ and R⁵ are independently selected from H and C₁-C₃ alkyl;
Z is CH or N; Y is -O- or -C(R⁴)(R⁵)-; q is 1; R⁸ and R⁹ are each H; Q is a substituted or
unsubstituted phenyl or furanyl group or a benzo[1,3]dioxyl or benzo[1,4]dioxyl group, where
the substituted phenyl or furanyl group contains one, two or three substituents selected from
halo, C₁-C₄ alkyl; C₁-C₄ alkylthio; or -NR^{Q1}R^{Q2}, where R^{Q1} and R^{Q2} taken together with the
nitrogen to which they are attached form a 4-7 membered heterocyclic ring, which may
optionally contain one or more additional heteroatoms selected from N, O and S; t is 0 or 1;
m is 0 or 1; R⁶ and R⁷ are independently selected from H and C₁-C₄ alkyl; W¹ is [~~methy~~l,]
unsubstituted phenyl, naphthyl, pyridyl, thienyl or pyrrolyl or substituted phenyl or pyridyl
containing one or two substituents independently selected from halo, alkyl and alkoxy,
specifically, chloro, methyl and methoxy; W² is C₁-C₄ alkyl, C₂-C₄ alkynyl, C₃-C₆ cycloalkyl,
aryl, Het hydroxy, aryloxy-, C₁-C₄ alkoxy-, -OCOC₁-C₄ alkyl, -OCOaryl, or -NR^{W1}R^{W2},
where R^{W1} and R^{W2} are independently H or C₁-C₄ alkyl or taken together with the nitrogen to
which they are attached form a 4-7 membered heterocyclic ring, which may optionally contain
one or more additional heteroatoms selected from N, O and S; W³ is H or C₁-C₄ alkyl; or a
pharmaceutically acceptable salt or [~~solvate~~] hydrate thereof.

67. (Currently amended, Withdrawn): The compound according to claim 49, wherein
X is chloro, bromo, cyano, carboxy-, methylcarboxy-, hydroxy, methoxy, methyl,
trifluoromethyl, 1,3-dihydroxy-prop-2-yl (-CH(CH₂OH)₂), isopropyl, n-butyl, isobutyl,
2,2-dimethylpropyl, phenylcarbonyl, triazolyl, tetrazolyl, -NH₂, -NHCH₃, -NHCH₂CH₃,
-NHCH₂CH₂CH₃, -NHCH₂CH₂CH₂CH₃, -NHCH₂CH₂CH₂CH₂CH₃, -NHCH₂C(CH₃)₃,
-NHCH₂CH(CH₃)₂, -NHCH₂CH₂CH(CH₃)₂, -NH-cyclopentyl, -NH-phenyl,
-NHCH₂-cyclopropyl, -NHCH(CH₃)₂, -NHCH₂CF₃, -N(CH₃)₂, -N(CH₂CH₃)₂,
-NHCH(CH₂CH₃)₂, -NHCH₂CH(CH₂CH₃)₂, -NHCH₂CH₂OH, -NHCH₂CO₂H,
-N(CH₃)CH₂CO₂H, -NHC(CH₃)₂CO₂H, -NHCH(CH₃)CO₂H, -(R)-NHCH(CH₃)CO₂H,

- (S)-NHCH(CH₃)CO₂H, -NHCH₂-1H-imidazol-2-yl, -NHCH₂-(1-CH₃-imidazol-2-yl, -NH-(pyrimidin-2-yl), -morpholin-4-yl, -thiomorpholin-4-yl, -piperidin-1-yl, -piperidin-1-yl-(4-carboxylic acid), -piperidin-1-yl-(4-acetic acid), -piperidin-4-yl-(1-acetic acid), -2,5-dimethyl-pyrrol-1-yl, -pyrrolidin-1-yl, -((R)-2-CO₂H-pyrrolidin-1-yl), -((S)-2-CO₂H-pyrrolidin-1-yl), -piperazin-1-yl, -(4-methyl-piperazin-1-yl), -piperazin-1-yl-(4-acetic acid), -NHCH₂-(5-bromo-thien-2-yl), -NHCH₂-1H-imidazol-2-yl, -NHCH₂-(1-methyl-imidazol-2-yl, -NHCOCH₃, -N(CH₃)COCH₃, -NHCO₂C(CH₃)₃, -NHCOCH₂CH₃, -NHCOC(CH₃)₂, -NHCO-furan-2-yl, -N(CH₃)CO-furan-2-yl, -NHCO-thien-2-yl, -NHCO-cyclopropyl, -NHCO-(5-bromo-thien-2-yl, -NHCO-(2,5-dimethyl-pyrrol-3-yl), -NHSO₂CH₃, -N(CH₃)SO₂CH₃, -NHSO₂CF₃, -NHSO₂phenyl, -N(CH₃)SO₂phenyl, -NHSO₂CH₂CH₃, -NHSO₂CH₂CF₃, -NHSO₂CH₂CH₂CH₃, -NHSO₂CH(CH₃)₂, -NHCONH(2-chlorophenyl), -N(CH₃)CONH(3,5-dimethoxyphenyl), -N(CH₃)CONH(2-chlorophenyl), -N(CH₃)CO-(benzo[1,3]diox-5-yl), -SO₂NHCH₃, and -SO₂N(CH₃)₂; p is 0, 1 or 2; R¹ and R² are H C₁-C₄ alkyl or R¹ and R² together with the carbon to which they are attached form a 3, 4 or 5 membered carbocyclic ring; Z is CH or N; k is 0 or k is 1 and R³ is methyl, trifluoromethyl, chloro or methoxy; n is 3 and R⁴ and R⁵ are independently selected from H and methyl; Y is -O- or -C(R⁴)(R⁵)-; q is 1; R⁸ and R⁹ are each H; Q is 2-chloro-3-(trifluoromethyl)phenyl, 3-methyl-4-fluoro-phenyl, 4-tert-butyl-phenyl, 4-(methylthio)phenyl, 2,4,5-trifluoro-phenyl, 4-isopropyl-phenyl, 5-(piperidin-1-yl)-furan-2-yl, benzo[1,3]diox-5-yl, or 2,3-dihydrobenzo[1,4]dioxin-6-yl; t is 0 or 1; m is 0 or 1; R⁶ and R⁷ are independently selected from H and methyl; W¹ is [~~[[methyl],~~] phenyl, naphth-1-yl, pyrid-2-yl, 4-methyl-pyrid-2-yl, thien-2-yl, thien-3-yl, pyrrol-2-yl, 2-chlorophenyl, 3-chlorophenyl, 4-chlorophenyl, 2-methoxyphenyl, or 4-methoxyphenyl; W² is methyl, ethyl, ethynyl, isopropyl, n-butyl, 2-methylpropyl, trifluoromethyl, cyclohexyl, unsubstituted phenyl, hydroxy, methoxy, phenoxy, dimethylamino, morpholin-4-yl, phenylcarbonyloxy, or methylcarbonyloxy; W³ is H or methyl; or a pharmaceutically acceptable salt or [~~[[solvate]]~~] hydrate thereof.

Claims 68-69 (Cancelled).

70. (Withdrawn): The compound according to claim 49, wherein W^1 and W^2 are not each independently C_3 - C_8 cycloalkyl or aryl or W^3 is not H or any one of R^6 or R^7 is not H or R^8 and R^9 are each C_1 - C_4 alkyl when:

X is $COOR^{10}$;

Z is CH or CR^3 and k is 0-4 or Z is N and k is 0-3;

p is 0-8;

n is 3;

q is 0 or 1;

Q is selected from optionally unsubstituted or substituted C_3 - C_8 cycloalkyl, phenyl and monocyclic Het;

each R^1 and R^2 is independently selected from H, C_1 - C_6 alkyl, -OH, -O- C_1 - C_6 alkyl, -SH, and -S- C_1 - C_6 alkyl; and

each R^3 is the same or different and is independently selected from halo, cyano, nitro, -CONR¹²R¹³, -COR¹⁴, -SR¹¹, -SO₂R¹¹, -SOR¹⁴, -OCOR¹⁴ and optionally unsubstituted or substituted C_1 - C_6 alkyl, C_3 - C_6 alkenyl, -5-6 membered-Het, -C₀- C_6 alkyl-CO₂R¹¹, or -C₀- C_6 alkyl-NR¹²R¹³.

71. (Previously presented): A pharmaceutical composition comprising the compound according to claim 49 and a pharmaceutically acceptable carrier or diluent.

Claims 72-91 (Cancelled).

92. (Withdrawn): A method for the prevention or treatment of an LXR mediated disease or condition, wherein said disease or condition is selected from atherosclerosis and inflammation, comprising administering a therapeutically effective amount of the compound according to claim 49.

93. (Withdrawn): The method according to claim 92, comprising administering a therapeutically effective amount of the compound according to claim 52.

94. (Withdrawn): The method according to claim 92, comprising administering a therapeutically effective amount of the compound according to claim 55.

95. (Withdrawn): The method according to claim 92, comprising administering a therapeutically effective amount of the compound according to claim 56.

96. (Withdrawn): The method according to claim 92, comprising administering a therapeutically effective amount of the compound according to claim 57.

97. (Withdrawn): The method according to claim 92, comprising administering a therapeutically effective amount of the compound according to claim 60.

98. (Withdrawn): The method according to claim 92, comprising administering a therapeutically effective amount of the compound according to claim 63.

99. (Withdrawn): The method according to claim 92, comprising administering a therapeutically effective amount of the compound according to claim 64.

100. (Withdrawn): The method according to claim 92, comprising administering a therapeutically effective amount of the compound according to claim 66.

101. (Withdrawn): The method according to claim 92, comprising administering a therapeutically effective amount of the compound according to claim 67.

102. (Currently amended, Withdrawn): The method according to claim 92, comprising administering a compound selected from:

~~2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]propoxy}phenyl)-ethanol,~~

2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]propoxy}phenyl)-ethanol,

(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenylethyl)-{3-[3-(1,2,4-triazol-3-ylmethyl)-phenoxy]-propyl}-amine,

(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{3-[3-(1,2,3,4-tetrazol-5-ylmethyl)-phenoxy]-propyl}-amine,

(*S*)-(2-chloro-3-trifluoromethyl-benzyl)-(2-phenyl-propyl)-{3-[3-(1,2,3,4-tetrazol-3-ylmethyl)-phenoxy]-propyl}-amine,

(*R*)-(2-chloro-3-trifluoromethyl-benzyl)-(2-phenyl-propyl)-{3-[3-(1,2,3,4-tetrazol-3-ylmethyl)-phenoxy]-propyl}-amine,

(*S*)-2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2-phenyl-propyl)amino]propoxy}-phenyl)acetic acid,

(*R*)-2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2-phenyl-propyl)amino]propoxy}-phenyl)acetic acid,

2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2-acetoxy-2-phenyl-ethyl)amino]propoxy}-phenyl)acetic acid,

(3-{3-[(2-acetoxy-2-phenyl-ethyl)-(2-chloro-3-trifluoromethyl-benzyl)-amino]-propoxy}-phenyl)-acetic acid methyl ester,

(3-{4-[(2-chloro-3-(trifluoromethyl)benzyl)-(2,2-diphenylethyl)-amino]butyl}phenyl)-acetic acid,

1-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-cyclobutanecarboxylic acid,

N-(2,2-diphenylethyl)-*N*-(2-chloro-3-trifluoromethylbenzyl)-3-(3-[2-hydroxy-2-methylpropyl]phenoxy)propylamine,

(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-{2-[(1*H*-imidazol-2-ylmethyl)-amino]-ethyl}-phenoxy)-propyl]-amine,

N-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-methanesulfonamide,

N-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-*N*-methyl-amine,

[2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethylamino]-acetic acid,

(*R*)-1-[2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethyl]-pyrrolidine-2-carboxylic acid,

furan-2-carboxylic acid *N*-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-amide,

N-(2-[3-chlorophenyl]-propyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine,
(2-chloro-3-trifluoromethyl-benzyl)-{3-[3-(2-morpholin-4-yl-ethyl)-phenoxy]-propyl}-((S)-2-phenyl-propyl amine,
[4-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-piperazin-1-yl]-acetic acid,
2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-((S)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-2-methyl-propionic acid,
(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{3-[3-(4-methyl-piperazin-1-yl)-phenoxy]-propyl}-amine,
(3-{(R)-3-[(2-chloro-3-trifluoromethyl-benzyl)-((S)-2-phenyl-propyl)-amino]-butoxy}-phenyl)-acetic acid,
[1-(3-{3-[(2-chloro-3-(trifluoromethyl)-benzyl)-(2,2-diphenylethyl)-amino]-propoxy}-phenyl)-piperidine-4-carboxylic acid,
[4-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-((S)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-piperazin-1-yl]-acetic acid,
[4-(3-{(R)-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-methyl-propoxy}-phenyl)-piperazin-1-yl]-acetic acid,
and a pharmaceutically acceptable salt or hydrate thereof.

103. (Withdrawn): A method for increasing reverse cholesterol transport, said method comprising administering a therapeutically effective amount of the compound according to claim 49.

104. (Withdrawn): A method for inhibiting cholesterol absorption, said method comprising administering a therapeutically effective amount of the compound according to claim 49.

105. (Currently amended): A compound selected from:

~~2-(3-{3-[[2-Chloro-3-(trifluoromethyl)benzyl](2,2-~~
~~diphenylethyl)amino]propoxy}phenyl)-ethanol~~ 2-(3-{3-[[2-Chloro-3-
(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]propoxy}phenyl)-ethanol; 2-(3-{3-[[2-

Chloro-3-(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]propoxy}-phenyl)acetic acid, N-oxide; (3-{3-[[2-Chloro-3-(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]propoxy}-bromobenzene; (4-{3-[[2-Chloro-3-(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]propoxy}-bromobenzene; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenylethyl)-{3-[3-(1,2,4-triazol-3-ylmethyl)-phenoxy]-propyl}-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{3-[3-(1,2,3,4-tetrazol-5-ylmethyl)-phenoxy]-propyl}-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2-cyclohexyl-2-phenyl-ethyl)-{3-[3-(1,2,3,4-tetrazol-5-ylmethyl)-phenoxy]-propyl}-amine; (S)-(2-Chloro-3-trifluoromethyl-benzyl)-(2-phenyl-propyl)-{3-[3-(1,2,3,4-tetrazol-3-ylmethyl)-phenoxy]-propyl}-amine; (R)-(2-Chloro-3-trifluoromethyl-benzyl)-(2-phenyl-propyl)-{3-[3-(1,2,3,4-tetrazol-3-ylmethyl)-phenoxy]-propyl}-amine; (S)-2-(3-{3-[[2-Chloro-3-(trifluoromethyl)benzyl](2-phenyl-propyl)amino]propoxy}-phenyl)acetic acid; (R)-2-(3-{3-[[2-Chloro-3-(trifluoromethyl)benzyl](2-phenyl-propyl)amino]propoxy}-phenyl)acetic acid; 2-(3-{3-[[2-Chloro-3-(trifluoromethyl)benzyl] naphthalen-1-ylmethyl-amino]propoxy}-phenyl)acetic acid; 2-(3-{3-[[2-Chloro-3-(trifluoromethyl)benzyl]-benzylamino]propoxy}-phenyl)acetic acid; 2-(3-{3-[[2-Chloro-3-(trifluoromethyl)-benzyl]phenethylamino]propoxy}-phenyl)acetic acid; 2-(3-{3-[[2-Chloro-3-(trifluoromethyl)benzyl](2-hydroxy-2-phenyl-ethyl)amino]propoxy}-phenyl)acetic acid; 2-(3-{3-[[2-Chloro-3-(trifluoromethyl)benzyl](2-acetoxy-2-phenyl-ethyl)amino]propoxy}-phenyl)acetic acid; 2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2-phenoxy-2-phenyl-ethyl)amino]propoxy}-phenyl)acetic acid; Benzoic acid 2-[3-(3-carboxymethyl-phenoxy){2-chloro-3-(trifluoromethyl)benzyl}propylamino]-1-phenyl ethyl ester; (3-{3-[(2-Acetoxy-2-phenyl-ethyl)-(2-chloro-3-trifluoromethyl-benzyl)-amino]-propoxy}-phenyl)-acetic acid methyl ester; Benzoic acid 2-[3-(3-methoxycarbonylmethyl-phenoxy){2-chloro-3-(trifluoromethyl)benzyl}propylamino]-1-phenyl ethyl ester; (3-{4-[(2-Chloro-3-(trifluoromethyl)benzyl)-(2,2-diphenylethyl)-amino]butyl}phenyl)-acetic acid; (3-{3-[(4-Fluoro-3-methyl-benzyl)-((R)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-acetic acid; (3-{3-[Benzo[1,3]dioxol-5-ylmethyl-((R)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-acetic acid; (3-{3-[(4-tert-Butyl-benzyl)-((R)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-acetic acid; (3-{3-[(2,3-Dihydro-benzo[1,4]dioxin-6-ylmethyl)-((R)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-acetic acid; (3-{3-[(4-Methylsulfanyl-benzyl)-((R)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-acetic acid; (3-{3-[(R)-2-Phenyl-propyl)-(2,4,5-trifluoro-benzyl)-amino]-

propoxy}-phenyl)-acetic acid; (3-{3-[(*(R)*)-2-Phenyl-propyl)-(5-piperidin-1-yl-furan-2-ylmethyl)-amino]-propoxy}-phenyl)-acetic acid; (3-{3-[(4-Isopropyl-benzyl)-(*(R)*)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-acetic acid; 2-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-propane-1,3-diol; *N*-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-carbamic acid tert-butyl ester; 3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethylamino]-propoxy}-phenylamine; *N*-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-acetamide; Furan-2-carboxylic acid *N*-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-amide; *N*-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-methanesulfonamide; *N*-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-benzenesulfonamide; 1-(2-Chloro-phenyl)-3-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-urea; *N*-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-*N*-methyl-amine; *N*-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-*N*-methyl-acetamide; Furan-2-carboxylic acid *N*-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-*N*-methyl-amide; *N*-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-*N*-methyl-methanesulfonamide; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-*N*-methyl-benzenesulfonamide; 3-(2-Chloro-phenyl)-1-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-1-methyl-urea; Benzo[1,3]dioxole-5-carboxylic acid *N*-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-*N*-methyl-amide; 1-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-3-(3,5-dimethoxy-phenyl)-1-methyl-urea; Propane-1-sulfonic acid (5-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-2-methyl-phenyl)-amide; 3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-2-methyl-phenylamine; 2-Chloro-5-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenylamine; 3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-phenyl)-cyclopentyl-amine; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-phenyl)-isopropyl-amine; Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-phenyl)-ethyl-amine; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-phenyl)-(3-methyl-butyl)-

amine; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-isobutyl-amine; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-phenyl)-(2,2,2-trifluoroethyl)-amine; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-phenyl)-cyclopropylmethyl-amine; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-phenyl)-(2-ethyl-butyl)-amine; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-phenyl)-(2,2-dimethyl-propyl)-amine; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-phenyl)-hexyl-amine; Butyl-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-phenyl)-amine; [1-(3-{3-[(2-Chloro-3-(trifluoromethyl)-benzyl)-(2,2-diphenylethyl)-amino]-propoxy}-phenyl)-piperidine-4-carboxylic acid; [1-(3-{3-[(2-Chloro-3-(trifluoromethyl)-benzyl)-(2,2-diphenylethyl)-amino]-propoxy}-phenyl)-piperidine-4-yl)-acetic acid; [4-(3-{3-[(2-Chloro-3-(trifluoromethyl)-benzyl)-(2,2-diphenylethyl)-amino]-propoxy}-phenyl)-piperidin-1-yl]-acetic acid; *rac*-±-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(trifluoro-phenyl-propyl)-amino]-propoxy}-phenyl)-acetic acid; *rac*-±-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2-dimethylamino-2-phenyl-ethyl)-amino]-propoxy}-phenyl)-acetic acid; *rac*-±-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2-morpholin-4-yl-2-phenyl-ethyl)-amino]-propoxy}-phenyl)-acetic acid; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(6-morpholin-4-yl-pyridin-2-yloxy)-propyl]-amine; [3-(6-Chloro-pyridin-2-yloxy)-propyl]-(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(6-(4-methyl-piperazin-1-yl)-pyridin-2-yloxy)-propyl]-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(6-piperazin-1-yl-pyridin-2-yloxy)-propyl]-amine; [4-(6-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-pyridin-2-yl)-piperazin-1-yl]-acetic acid; 2-(3-{3-[(2-Chloro-3-(trifluoromethyl)benzyl)](S)-2-phenyl-propyl)amino)-(R)-1-methyl-propoxy}-phenyl)acetic acid; 2-(3-{3-[(2-Chloro-3-(trifluoromethyl)benzyl)](S)-2-phenyl-propyl)amino)-(R)-1-methyl-propoxy}-phenyl)ethanol; 2-(3-{3-[(2-Chloro-3-(trifluoromethyl)benzyl)](S)-2-phenyl-propyl)amino)-(R)-2-methyl-propoxy}-phenyl)acetic acid; 2-(3-{3-[(2-Chloro-3-(trifluoromethyl)benzyl)](S)-2-phenyl-propyl)amino)-(R)-2-methyl-propoxy}-phenyl)ethanol; 2-(3-{3-[(2-Chloro-3-(trifluoromethyl)benzyl)](R)-2-phenyl-propyl)amino)-(R)-2-methyl-propoxy}-phenyl)acetic acid; 2-(3-{3-[(2-Chloro-3-(trifluoromethyl)benzyl)](R)-2-phenyl-propyl)amino)-(R)-2-methyl-propoxy}-phenyl)ethanol; (R)-2-(3-{3-[(2-Chloro-3-(trifluoromethyl)benzyl)](2,2-

diphenylethyl)amino]-2-methyl-propoxy)-phenyl)ethanol; 3-{3-[(3-Chloro-2-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy-*N,N*-dimethyl-benzenesulfonamide; Cyclopropanecarboxylic acid 3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-benzylamide; N-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-benzyl)-isobutyramide; Acetic acid (3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-benzylcarbamoyl)-methyl ester; N-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-benzyl)-propionamide; 2,5-Dimethyl-2-H-pyrazole-3-carboxylic acid 3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-benzylamide; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-(3-*o*-tolylloxy-propyl)-amine; 2-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-benzonitrile; 3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-benzonitrile; [3-(3-Chloro-phenoxy)-propyl]-(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(2-methoxy-phenoxy)-propyl]-amine; [3-(2-Chloro-phenoxy)-propyl]-(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-(3-phenoxy-propyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-isopropyl-phenoxy)-propyl]-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(4-methoxy-phenoxy)-propyl]-amine; 3-{3-[(Chloro-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenol; 2-{3-[(Chloro-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenol; 3-{3-[(Chloro-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenylamine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-trifluoromethyl-phenoxy)-propyl]-amine; 1-(3-{3-[(Chloro-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethanone; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-phenylamine; [3-(Benzo[1,3]dioxol-5-yloxy)-propyl]-(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-(3-*m*-tolylloxy-propyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-methoxy-phenoxy)-propyl]-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-isobutyl-phenoxy)-propyl]-amine; [3-(3-Butyl-phenoxy)-propyl]-(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-{3-[3-(2,2-dimethyl-propyl)-phenoxy]-propyl}-(2,2-diphenyl-ethyl)-amine; (4-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-benzyl)-methyl-amine; (2-Chloro-3-

trifluoromethyl-benzyl)-[3-(4-dimethylaminomethyl-phenoxy)-propyl]-(2,2-diphenyl-ethyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(4-morpholin-4-ylmethyl-phenoxy)-propyl]-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-[4-(4-methyl-piperazin-1-ylmethyl)-phenoxy]-propyl]-amine; (3-{3-[(Chloro-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-benzyl)-methyl-amine; (2-Chloro-3-trifluoromethyl-benzyl)-[3-(3-dimethylaminomethyl-phenoxy)-propyl]-(2,2-diphenyl-ethyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-morpholin-4-ylmethyl-phenoxy)-propyl]-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-[3-(4-methyl-piperazin-1-ylmethyl)-phenoxy]-propyl]-amine; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-benzyl)-isopropyl-amine; {3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-4-trifluoromethyl-phenylamine; {3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-4-methyl-phenylamine; Ethanesulfonic acid (3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-4-methyl-phenyl)-amide; Propane-2-sulfonic acid (3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-4-methyl-phenyl)-amide; Methanesulfonic acid (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-4-methyl-phenyl)-amide; 2,2,2-Trifluoro-ethanesulfonic acid (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-4-methyl-phenyl)-amide; Ethanesulfonic acid (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-phenyl)-amide; 2,2,2-Trifluoro-ethanesulfonic acid (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-phenyl)-amide; *N*-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-phenyl)-1,1,1-trifluoro-methanesulfonamide; Propane-2-sulfonic acid (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-phenyl)-amide; {3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-4-methoxy-phenylamine; Ethanesulfonic acid (3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl-amino)]-propoxy}-4-methoxy-phenyl)-amide; (2-Chloro-3-trifluoromethyl-benzyl)-[3-[3-(2-morpholin-4-yl-ethyl)-phenoxy]-propyl]-((*S*)-2-phenyl-propyl amine; (2-Chloro-3-trifluoromethyl-benzyl)-[3-[3-(2-ethylamino-ethyl)-phenoxy]-propyl]-((*S*)-2-phenyl-propyl)-amine; (3-{(*R*)-3-[(2-Chloro-3-trifluoromethyl-benzyl)-((*S*)-2-phenyl-propyl)-amino]-butoxy}-phenyl)-acetic acid; (3-{(*S*)-3-[(2-Chloro-3-trifluoromethyl-benzyl)-((*S*)-2-phenyl-propyl)-amino]-butoxy}-phenyl)-acetic acid; 2-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-

((S)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-ethanol; 2-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-((S)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-2-methyl-propionic acid; 2-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-((R)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-2-methyl-propionic acid; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2-thiophen-3-yl-propyl)-amino]-propoxy}-phenyl)-acetic acid; 2-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2-thiophen-3-yl-propyl)-amino]-propoxy}-phenyl)-ethanol; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2-thiophen-2-yl-propyl)-amino]-propoxy}-phenyl)-acetic acid; (3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2-pyridin-2-yl-propyl)-amino]-propoxy}-phenyl)-acetic acid; [3-(3-{(2-Chloro-3-trifluoromethyl-benzyl)-[2-(4-methyl-pyridin-2-yl)-propyl]-amino}-propoxy)-phenyl]-acetic acid; [3-(3-{(2-Chloro-3-trifluoromethyl-benzyl)-[3,3,3-trifluoro-2-(1H-pyrrol-2-yl)-propyl]-amino}-propoxy)-phenyl]-acetic acid; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-(3-{3-[2-(4-methyl-piperazin-1-yl)-ethyl]-phenoxy}-propyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{3-[3-(2-methylamino-ethyl)-phenoxy]-propyl}-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-{2-[(1H-imidazol-2-ylmethyl)-amino]-ethyl}-phenoxy)-propyl]-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{3-[3-(2-ethylamino-ethyl)-phenoxy]-propyl}-amine; [3-(3-{2-[(5-Bromo-thiophen-2-ylmethyl)-amino]-ethyl}-phenoxy)-propyl]-[2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-{2-[(thiophen-2-ylmethyl)-amino]-ethyl}-phenoxy)-propyl]-amine; (2-Chloro-3-trifluoromethyl-benzyl)-{3-[3-(2-dimethylamino-ethyl)-phenoxy]-propyl)-(2,2-diphenyl-ethyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{3-[3-(2-pyrrolidin-1-yl-ethyl)-phenoxy]-propyl}-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{3-[3-(2-morpholin-4-yl-ethyl)-phenoxy]-propyl}-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{(R)-1-methyl-3-[3-(2-morpholin-4-yl-ethyl)-phenoxy]-propyl}-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{(R)-2-methyl-3-[3-(2-morpholin-4-yl-ethyl)-phenoxy]-propyl}-amine; {3-[3-(2-Amino-ethyl)-phenoxy]-propyl}-(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amine; [2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethyl]-isopropyl-amine; [2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethyl]-propyl-amine; 2-[2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethylamino]-ethanol; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-{2-[(1-methyl-1H-imidazol-2-ylmethyl)-

amino]-ethyl)-phenoxy)-propyl]-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{3-[3-(2-thiomorpholin-4-yl-ethyl)-phenoxy]-propyl]-amine; [2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethylamino]-acetic acid; [2-(3-{(R)-3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-butoxy}-phenyl)-ethylamino]-acetic acid; {[2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethyl]-methyl-amino}-acetic acid; 2-[2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethylamino]-2-methyl-propionic acid; (S)-2-[2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethylamino]-propionic acid; (R)-1-[2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethyl]-pyrrolidine-2-carboxylic acid; (S)-1-[2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethyl]-pyrrolidine-2-carboxylic acid; [2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethyl]-pyrimidin-2-yl-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-morpholin-4-yl-phenoxy)-propyl]-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-piperidin-1-yl-phenoxy)-propyl]-amine; (3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-diethyl-amine; (2-Chloro-3-trifluoromethyl-benzyl)-{3-[3-(2,5-dimethyl-pyrrol-1-yl)-phenoxy]-propyl}-(2,2-diphenyl-ethyl)-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-piperazin-1-yl-phenoxy)-propyl]-amine; (2-Chloro-3-trifluoromethyl-benzyl)-((S)-2-phenyl-propyl)-[3-(3-piperazin-1-yl-phenoxy)-propyl]-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[(R)-2-methyl-3-(3-piperazin-1-yl-phenoxy)-propyl]-amine; ~~(2-Chloro-3-trifluoromethyl-benzyl)-isobutyl-[3-(3-piperazin-1-yl-phenoxy)-propyl]-amine~~; [4-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-piperazin-1-yl]-acetic acid; [4-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-methyl-propoxy}-phenyl)-piperazin-1-yl]-acetic acid; [4-(3-{(R)-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-methyl-propoxy}-phenyl)-piperazin-1-yl]-acetic acid; ~~[4-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-isobutyl-amino]-propoxy}-phenyl)-piperazin-1-yl]-acetic acid~~; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{3-[3-(4-methyl-piperazin-1-yl)-phenoxy]-propyl}-amine; (2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-pyrrolidin-1-yl-phenoxy)-propyl]-amine; (3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenylamino)-acetic acid; [(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-N-methyl-

amino]-acetic acid; N-(2,2-Diphenylethyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-[3-(2-methyl-2-aminopropyl)phenoxy]propylamine; N-(2,2-Diphenylethyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-[2-hydroxymethyl]phenoxy)propylamine; N-(2,2-Diphenylethyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-[2-hydroxy-2-methylpropyl]phenoxy)propylamine; N-(2,2-Diphenylethyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-N-methylsulfonamidophenoxy)propylamine; N-(2-[2-Chlorophenyl]-propyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; N-(2-[3-Chlorophenyl]-propyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; N-(2-[4-Chlorophenyl]-propyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; N-(2-[2-Methoxyphenyl]-propyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; N-(2-[4-Methoxyphenyl]-propyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; N-(2-Phenyl-4-methylpentyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; N-(2-Phenylbutyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; N-(2-[2-Methyl-2-phenyl]propyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; N-(2-Phenyl-3-methylbutyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; N-(2-Phenylhexyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; N-(2-Phenyl-3-butynyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; (S)-N-(2-Phenyl-2-methoxyethyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; (R)-N-(2-Phenyl-2-methoxyethyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine; (R)-N-(2-Phenyl-2-methoxyethyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-[2-hydroxy-2-methylpropyl]phenoxy)propylamine; 2-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)acetic acid; 1-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-cyclobutanecarboxylic acid; 1-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-cyclopentanecarboxylic acid; 1-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-amino]-propoxy}-phenyl)-cyclopropanecarboxylic acid;

and a pharmaceutically acceptable salt or hydrate thereof.

106. (Currently amended, Withdrawn): The compound according to claim 49, selected from:

~~2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]propoxy}phenyl)-ethanol~~

2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]propoxy}phenyl)-ethanol,

(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenylethyl)-{3-[3-(1,2,4-triazol-3-ylmethyl)-phenoxy]-propyl}-amine,

(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{3-[3-(1,2,3,4-tetrazol-5-ylmethyl)-phenoxy]-propyl}-amine,

(2-chloro-3-trifluoromethyl-benzyl)-(2-cyclohexyl-2-phenyl-ethyl)-{3-[3-(1,2,3,4-tetrazol-5-ylmethyl)-phenoxy]-propyl}-amine,

2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl]-benzylamino]propoxy}-phenyl)acetic acid,

2-(3-{3-[[2-chloro-3-(trifluoromethyl)-benzyl]phenethylamino]propoxy}-phenyl)acetic acid,

2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2-hydroxy-2-phenyl-ethyl)amino]propoxy}-phenyl)acetic acid,

2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2-acetoxy-2-phenyl-ethyl)amino]propoxy}-phenyl)acetic acid,

2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2-phenoxy-2-phenyl-ethyl)amino]propoxy}-phenyl)acetic acid,

(3-{3-[(2-acetoxy-2-phenyl-ethyl)-(2-chloro-3-trifluoromethyl-benzyl)-amino]-propoxy}-phenyl)-acetic acid methyl ester,

benzoic acid 2-[3-(3-methoxycarbonylmethyl-phenoxy){2-chloro-3-(trifluoromethyl)benzyl}propylamino]-1-phenyl ethyl ester,

(3-{4-[(2-chloro-3-(trifluoromethyl)benzyl)-(2,2-diphenylethyl)-amino]butyl}phenyl)-acetic acid,

~~2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]propoxy}phenyl)-ethanol,~~

~~(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenylethyl)-{3-[3-(1,2,4-triazol-3-ylmethyl)-phenoxy]-propyl}-amine,~~

~~(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-[3-(1,2,3,4-tetrazol-5-ylmethyl)-phenoxy]-propyl]-amine,~~

~~furan-2-carboxylic acid N-(3-{3-[(2-Chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-amide,~~

~~2-(3-{3-[(2-chloro-3-(trifluoromethyl)benzyl)](2-acetoxy-2-phenyl-ethyl)amino}propoxy)-phenyl)acetic acid,~~

~~(3-{3-[(2-acetoxy-2-phenyl-ethyl)-(2-chloro-3-trifluoromethyl-benzyl)-amino]-propoxy}-phenyl)-acetic acid methyl ester,~~

~~(3-{4-[(2-chloro-3-(trifluoromethyl)benzyl)-(2,2-diphenylethyl)-amino]butyl}phenyl)-acetic acid,~~

~~1-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-cyclobutanecarboxylic acid,~~

~~N-(2,2-diphenylethyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-[2-hydroxy-2-methylpropyl]phenoxy)propylamine,~~

~~(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(3-{2-[(1H-imidazol-2-ylmethyl)-amino]-ethyl}-phenoxy)-propyl]-amine,~~

~~N-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-methanesulfonamide,~~

~~N-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-2,2-diphenylethyl-amino]-propoxy}-phenyl)-N-methyl-amine,~~

~~[2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethylamino]-acetic acid,~~

~~N-(2-[3-chlorophenyl]-propyl)-N-(2-chloro-3-trifluoromethylbenzyl)-3-(3-carboxymethylenephenoxy)propylamine,~~

~~[4-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-piperazin-1-yl]-acetic acid,~~

~~(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-{3-[3-(4-methyl-piperazin-1-yl)-phenoxy]-propyl}-amine,~~

~~[1-(3-{3-[(2-chloro-3-(trifluoromethyl)-benzyl)-(2,2-diphenylethyl)-amino]-propoxy}-phenyl)-piperidine-4-carboxylic acid,~~

~~and a pharmaceutically acceptable salt or hydrate thereof.~~

107. (Withdrawn): The compound according to claim 49, selected from:

(S)-(2-chloro-3-trifluoromethyl-benzyl)-(2-phenyl-propyl)-{3-[3-(1,2,3,4-tetrazol-3-ylmethyl)-phenoxy]-propyl}-amine,

(R)-(2-chloro-3-trifluoromethyl-benzyl)-(2-phenyl-propyl)-{3-[3-(1,2,3,4-tetrazol-3-ylmethyl)-phenoxy]-propyl}-amine,

(S)-2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2-phenyl-propyl)amino]propoxy}-phenyl)acetic acid,

(R)-2-(3-{3-[[2-chloro-3-(trifluoromethyl)benzyl](2-phenyl-propyl)amino]propoxy}-phenyl)acetic acid,

(R)-1-[2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-propoxy}-phenyl)-ethyl]-pyrrolidine-2-carboxylic acid,

(2-chloro-3-trifluoromethyl-benzyl)-{3-[3-(2-morpholin-4-yl-ethyl)-phenoxy]-propyl}-((S)-2-phenyl-propyl amine,

2-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-((S)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-2-methyl-propionic acid,

(3-{(R)-3-[(2-chloro-3-trifluoromethyl-benzyl)-((S)-2-phenyl-propyl)-amino]-butoxy}-phenyl)-acetic acid,

[4-(3-{3-[(2-chloro-3-trifluoromethyl-benzyl)-((S)-2-phenyl-propyl)-amino]-propoxy}-phenyl)-piperazin-1-yl]-acetic acid,

[4-(3-{(R)-[(2-chloro-3-trifluoromethyl-benzyl)-diphenylethyl-amino]-methyl-propoxy}-phenyl)-piperazin-1-yl]-acetic acid,

and a pharmaceutically acceptable salt or hydrate thereof.